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CENTRAL FAX CENTER

Amendments to the Claims:

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This listing of claims will replace all prior versions and listings of claims in this application.

Listing of Claims:

- 1.-2. (Cancelled).
- (Previously Presented) A process for producing polyethylene using a loop reactor comprising:

using a mathematical model to predict a plurality of process control parameters based on desired product properties and reactor characteristics and controlling the process using the predicted process control parameters;

wherein variables used to prepare the mathematical model include a plurality of variables selected from the group consisting of: ethylene flow into reactor, ethylene flow out of reactor, isobutane flow into reactor, isobutane flow out of reactor, hexene flow into reactor, hexene flow out of reactor, hexene conversion in reactor, hydrogen flow into reactor, hydrogen flow out of reactor, hydrogen conversion in reactor, polymer flow out of reactor, liquid flow out of reactor, total mass flow into reactor, total mass flow out of reactor, total volume flow out of reactor, catalyst flow into reactor, catalyst flow out of reactor, concentration of ethylene in the reactor liquid, concentration of hexene in the reactor liquid, concentration of hydrogen in the reactor liquid, temperature of reactor, pressure of reactor, weight concentration of solids in a reactor slurry, volume concentration of solids in a reactor slurry, weight concentration of solids in [the] a settling leg solids bed, number of settling legs, reactor volume, settling leg diameter, settling leg height, bulk density of reactor polymer, density of reactor polymer, density of reactor[,] liquid, density of reactor slurry, residence time of reactor solids, catalyst activity, catalyst productivity, catalyst diameter, catalyst feed factor, catalyst activity factor, terminal velocity of settling polymer, polymer settling rate, rate of polymer leaving reactor that is not part of the settling

leg solids bed, rate of slurry leaving reactor that is part of the settling leg solids bed, rate of slurry leaving reactor that is not part of the settling leg solids bed, viscosity of reactor liquid, density of catalyst, polymer diameter, Archimedes number for polymer settling in settling leg, Reynolds number for polymer settling in settling leg, acceleration due to gravity, cross sectional area of a settling leg occupied by polymer, and cross sectional area of a settling leg; and

wherein the variables used to prepare the mathematical model are: polymer flow out of the reactor, liquid flow out of the reactor, concentration of ethylene in the reactor liquid, temperature of the reactor, pressure of the reactor, weight concentration of solids in the reactor slurry, reactor volume, settling leg diameter, settling leg height, bulk density of the reactor polymer, density of the reactor polymer, and density of the catalyst.

4. (Previously Presented) A process for producing polyethylene using a loop reactor comprising:

using a mathematical model to predict a plurality of process control parameters based on desired product properties and reactor characteristics and controlling the process using the predicted process control parameters;

wherein variables used to prepare the mathematical model include a plurality of variables selected from the group consisting of: ethylene flow into reactor, ethylene flow out of reactor, isobutane flow into reactor, isobutane flow out of reactor, hexene flow into reactor, hexene flow out of reactor, hexene conversion in reactor, hydrogen flow into reactor, hydrogen flow out of reactor, hydrogen conversion in reactor, polymer flow out of reactor, liquid flow out of reactor, total mass flow into reactor, total mass flow out of reactor, total volume flow out of reactor, catalyst flow into reactor, catalyst flow out of reactor, concentration of ethylene in the reactor liquid, concentration of hexene in the reactor liquid, concentration of hydrogen in the reactor liquid, temperature of reactor, pressure of reactor, weight concentration of solids in a reactor slurry, volume concentration of solids in a reactor slurry, weight concentration of solids in the settling leg solids bed, number of settling legs, reactor volume, settling leg diameter, settling leg height, bulk density of reactor polymer, density of reactor polymer, density of reactor[,] liquid, density of reactor slurry, residence time of reactor solids, catalyst activity, catalyst productivity, catalyst diameter, catalyst feed factor, catalyst activity factor, terminal velocity of settling polymer, polymer settling rate, rate of polymer leaving reactor that is not part of the settling leg solids bed, rate of slurry leaving reactor that is part of the settling leg solids bed, rate of slurry leaving reactor that is not part of the settling leg solids bed, viscosity of reactor liquid, density of catalyst, polymer diameter, Archimedes number for polymer settling in settling leg, Reynolds number for polymer settling in settling leg, acceleration due to gravity, cross sectional area of a settling leg occupied by polymer, and cross sectional area of a settling leg; and

wherein the variables used to prepare the mathematical model are: polymer flow out of the reactor, liquid flow out of the reactor, concentration of

ethylene in the reactor liquid, temperature of the reactor, pressure of the reactor, weight concentration of solids in the reactor slurry, reactor volume, settling leg diameter, settling leg height, bulk density of the reactor polymer, density of the reactor polymer, density of the catalyst, concentration of hexene in the reactor liquid, concentration of hydrogen in the reactor liquid, hexene conversion in the reactor, and hydrogen conversion in the reactor.

- (Cancelled).
- 6. (Cancelled).

- 7. (Cancelled).
- 8.-11. (Cancelled).
- 12. (Previously Presented) A process for designing a configuration of a loop reactor used to produce polymers comprising:

using a mathematical model to predict a plurality of process control parameters based on desired product properties and reactor characteristics and controlling the process using the predicted process control parameters, where variables used to prepare the mathematical model are selected from the group consisting of: mass balance of the reactor contents, reactor geometry, catalyst kinetics, and settling phenomena in [the] an outlet settling leg[s], which is used to determine relationships between inlet feed rates and reactor geometry on production conditions; and

wherein variability is decreased by 70% in comparison to a conventional PID controller.

- 13. (Previously Presented) The process of Claim 3, wherein the model is displayed in a spreadsheet.
- 14. (Previously Presented) The process of Claim 4, wherein the model is displayed in a spreadsheet.
- 15. (Previously Presented) The process of Claim 3, further comprising the steps of:

using sensors in the reactor to provide inputs to a PID controller; using another controller to receive data from said PID controller; and reprogramming the PID controller with the other controller, based upon the inputs from the reactor sensors.

- 16. (Previously Presented) The process of Claim 15, wherein a plurality of PID controllers are used to receive inputs from the reactor sensors, and wherein the PID controllers are reprogrammed based upon the inputs from the reactor sensors and the other PID controllers.
- 17. (Previously Presented) The process of Claim 3, wherein said mathematical model is incorporated into logic circuits of a controller.
- 18. (Previously Presented) The process of Claim 4, wherein said mathematical model is incorporated into logic circuits of a controller.
- 19. (Previously Presented) The process of Claim 3, wherein said mathematical model is created with a Fortran or C++ computer program.
- 20. (Previously Presented) The process of Claim 4, wherein said mathematical model is created with a Fortran or C++ computer program.
- 21. (Cancelled).

22. (Previously Presented) A process for producing polyethylene using a loop reactor comprising:

using a mathematical model to predict a plurality of process control parameters based on desired product properties and reactor characteristics and controlling the process using predicted and actual process control parameters;

wherein the mathematical model includes settling leg data and a plurality of variables selected from the group consisting of:

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Εi
                    ethylene flow into the reactor,
                    ethylene flow out of the reactor,
E<sub>o</sub>
                    isobutane flow into the reactor,
I_i
                    isobutane flow out of the reactor,
I_o
                    hexene flow into the reactor,
H_i
                    hexene flow out of the reactor,
H_o
                    hexene conversion in the reactor,
H_c
                    hydrogen flow into the reactor,
hi
                    hydrogen flow out of the reactor,
ho
              =
              =
                    hydrogen conversion in the reactor,
h_c
                    polymer flow out of the reactor,
P_o
                    liquid flow out of the reactor,
L_{\alpha}
              =
              =
                    total mass flow into the reactor,
T_{iw}
                   total mass flow out of the reactor,
T_{ow}
T_{ov}
              =
                    total volume flow out of the reactor,
                    catalyst flow into the reactor,
Ci
                    catalyst flow out of the reactor,
              =
C_{O}
                    concentration of ethylene in a reactor liquid,
C_{F}
              =
                    concentration of hexene in a reactor liquid,
C_H
C_h
                    concentration of hydrogen in a reactor liquid,
              =
                    temperature of the reactor.
T
              =
P
                    pressure of the reactor.
              =
S_w
              =
                    weight concentration of solids in a reactor slurry,
                    volume concentration of solids in a reactor slurry.
Sv
                    weight concentration of solids in a settling leg solids
S_{ws}
              =
                     bed.
N_t
              =
                    number of settling legs,
                    reactor volume.
dı
                    settling leg diameter,
                    settling leg height,
hı
              =
                    bulk density of a reactor polymer,
              =
r<sub>b</sub>
                    density of a reactor polymer,
rp
                    density of a reactor liquid,
              =
r
                    density of a reactor slurry,
rs
                    residence time of reactor solids,
              =
ts
k
                    catalyst activity,
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p	=	catalyst productivity,
d _c	=	catalyst diameter,
f_c	=	catalyst feed factor,
K _c	=	catalyst activity factor,
V_t	=	terminal velocity of a settling polymer,
$\dot{P_{\rm s}}$	=	polymer settling rate,
P_{n}	=	rate of a polymer leaving the reactor that is not part of
		the settling leg solids bed,
\mathcal{T}_{s}	=	rate of the slurry leaving the reactor that is part of the
		settling leg solids bed,
\mathcal{T}_{n}	=	rate of the slurry leaving the reactor that is not part of
		the settling leg solids bed,
m_t	=	viscosity of the reactor liquid,
r _c	=	density of the catalyst,
d_p	=	polymer diameter,
N _{Ar}	=	Archimedes number for polymer settling in a settling
		leg,
N_{Re}	=	Reynolds number for polymer settling in a settling leg,
\boldsymbol{g}	=	acceleration due to gravity,
Alp	=	cross sectional area of a settling leg occupied by
•		the polymer; and
A_{i}	=	cross sectional area of a settling leg.
,		

23. (Previously Presented) The process of Claim 22, wherein the mathematical model includes formulas selected from the group consisting of: mass balance of a reactor, determined by:

$$T_{bw} - (L_a + P_a) = 0$$
, where:

 P_o is polymer flow out of the reactor, L_o is liquid flow out of the reactor, and T_{iw} is total mass flow into the reactor;

mass flows in an outlet liquid based upon outlet concentrations and hydrogen flow, determined by:

$$E_o = \frac{L_o C_E}{100}$$
, $H_o = \frac{L_o C_H}{100}$, $h_o = \frac{L_o C_h}{100}$, $I_o + E_o + H_o = L_o$, and

 $I_o = L_o - E_o - H_o$, where:

 E_o is ethylene flow out of the reactor,

 L_o is liquid flow out of the reactor,

 C_E is concentration of ethylene in the reactor liquid,

 H_o is hexene flow out of the reactor,

 C_H is concentration of hexene in the reactor liquid.

 C_h is concentration of hydrogen in the reactor liquid, h_o is hydrogen flow out of the reactor, I_o is isobutane flow out of the reactor, and where h_o that is a smallest portion of total liquid flow;

an assumed mass balance of isobutane around the reactor, determined by:

$$I_i - I_o = 0$$
 , where:

 I_i is isobutane flow into the reactor, and I_0 is isobutane flow out of the reactor;

hexene and hydrogen feeds into the reactor, determined by:

$$H_o = \frac{H_i(100 - C_H)}{100}$$
, $H_i = \frac{100H_o}{100 - C_H}$, $h_o = \frac{h_i(100 - C_h)}{100}$, and

$$h_o = \frac{h_i (100 - C_h)}{100}$$
, where.

 H_l is hexene flow into the reactor,

 H_0 is hexene flow out of the reactor, and

 C_H is concentration of hexene in the reactor liquid, wherein a constant conversion of hexene and hydrogen into polymer across an intended operating region is assumed;

catalyst flow into the reactor, calculated using catalyst productivity and catalyst mass balance around the reactor, by:

$$p = \frac{P_o}{c_o}$$
 and $c_i - c_o = 0$, where:

p is catalyst productivity,

 P_o is polymer flow out of the reactor,

 c_i is catalyst flow into the reactor, and

 c_{o} is catalyst flow out of the reactor;

slurry density, calculated using liquid density, an assumed polymer density, and a solids concentration, determined by:

$$\rho_{s} = \frac{100\rho_{p}\rho_{l}}{100\rho_{p} + S_{w}(\rho_{l} - \rho_{p})};$$

volumetric flow rate of slurry out of the reactor determined from mass flow rate and density of each component, determined by:

$$T_{ov} = \frac{P_o}{\rho_p} + \frac{L_o}{\rho_l}$$

volumetric solids concentration in the reactor, determined by:

$$S_{\nu} = \frac{100}{1 + \frac{\rho_{p}}{\rho_{l}} \left(\frac{100}{S_{w}} - 1 \right)};$$

solid residence time, determined by:

$$\tau_s = \frac{V \rho_s S_W}{100 P_O};$$

catalyst activity, determined by:

$$p \approx \frac{P_o}{c_o}$$
 and $k = \frac{p}{\tau_s C_E}$;

a catalyst feed factor, determined by:

$$f_c = kc_i$$
;

a catalyst feed factor, using a mass balance of catalyst around the reactor, determined by:

$$f_c = kc_i = kc_o = \frac{P_o}{\tau_s C_E};$$

a catalyst activity factor that quantifies reactivity of a catalyst with respect to the catalyst size, by:

$$k_c = \frac{k}{d_c^3};$$

concentration of a polymer in a settling bed as a function of the polymer, liquid densities, and polymer bulk density, determined by:

$$S_{ws} = \frac{100}{1 + \rho_l \left(\frac{1}{\rho_b} - \frac{1}{\rho_p} \right)};$$

mass balance of two portions of an outlet slurry, determined by:

$$T_{ow} = T_s + T_n$$
;

mass balance of two portions of a polymer in outlet slurry, determined by:

$$P_o = P_s + P_n$$
;

an amount of polymer leaving a settling leg from the settled bed of polymer, calculated using known polymer concentrations of two polymer portions, by:

$$T_s = \frac{100P_s}{S_{ws}}$$
, $T_n = \frac{100P_n}{S_{ws}}$, $T_{ow} = \frac{100P_s}{S_{ws}} + \frac{100P_n}{S_{ws}}$,

$$T_{ow} = \frac{100P_s}{S_{ws}} + \frac{100(P_o - P_s)}{S_w}$$
, and $P_s = \frac{\frac{T_{ow}}{100} - \frac{P_o}{S_w}}{\frac{1}{S_{ws}} - \frac{1}{S_w}}$, where:

 T_s is the rate of slurry leaving the reactor that is part of the settling leg solids bed, T_n is the rate of slurry leaving reactor that is not part of the settling leg solids bed, T_{ow} is the total mass flow out of reactor, and wherein the flow rate indicates polymer settling rate in the settling leg;

viscosity of isobutane liquid as a function of temperature, determined by:

$$\mu_I(\text{cP}) = e^{-7.3891 + \frac{2582.6}{172.23 + T(\text{K})}}$$

polymer diameter, determined by:

$$d_p = 0.42 d_c \left(\frac{\rho_c p}{\rho_p} \right)^{1/3};$$

Archimedes number for polymer settling through the isobutane liquid, determined by:

$$N_{Ar} = \frac{d^3 \rho_l g(\rho_p - \rho_l)}{{\mu_l}^2}$$
 and $N_{Ar} = 1.13e05$;

calculation of a Reynolds number from the Archimedes number, by:

$$N_{\text{Re}} = \left(\sqrt{14.42 + 1.827\sqrt{N_{Ar}}} - 3.798\right)^2$$
, then

determining a velocity of a settling polymer, using:

$$N_{\text{Re}} = \frac{d_p v_t \rho_l}{\mu_l}$$
 and $v_t = \frac{N_{\text{Re}} \mu_l}{d_p \rho_l}$;

a mass flow of settling solids as related to a velocity of said solids, determined by:

$$P_s = N_i A_{l\rho} v_t \rho_{\rho};$$

an area of a settling leg occupied by polymer, determined by:

$$A_{lp} = \frac{A_l S_v}{100} = \frac{\pi}{400} d_l^2 S_v,$$

then substituting said area of the settling leg into the mass flow of settling solids, determined by:

$$P_{s} = \frac{\pi}{400} N_{l} d_{l}^{2} S_{v} v_{t} \rho_{p}$$
; and

wherein a number of settling legs required for reactor geometry and production conditions, is determined by:

$$N_l = \frac{400 P_s}{\pi d_l^2 S_v v_t \rho_p}.$$

24. (Previously Presented) The process of Claim 22, wherein it is assumed that HDPE is produced from monomers consisting of ethylene and hexene, and where ethylene feed into the reactor is calculated from a mass balance of ethylene around the reactor, using:

$$E_i - \{E_o + [P_o - (H_i - H_o)]\} = 0$$
 or $E_i = E_o + [P_o - (H_i - H_o)]$, where:

 E_i is ethylene flow into the reactor,

Eo is ethylene flow out of the reactor,

Po is polymer flow out of the reactor,

 H_i is hexene flow into the reactor, and

Ho is hexene flow out of the reactor.

25. (Previously Presented) The process of Claim 22, wherein it is assumed that HDPE is produced from monomers and where HDPE production kinetics are assumed to be first-order for ethylene concentration in a liquid and catalyst flow rate, wherein kinetic rate is determined by:

$$P_o = kc_o C_\kappa \tau_S$$

26. (Previously Presented) The process of Claim 23, where catalyst activity is determined by:

$$p = \frac{P_o}{c_o}$$
; and
$$k = \frac{p}{\tau_s C_E} = \frac{1624.9 \text{ lb/lb}}{(45.9 \text{ min})(5.0537 \text{ wt\%})} = 7.00 \text{ lb/lb/min/wt\%};$$

where catalyst feed factor is a function of other measured reactor variables, making it independent of catalyst type or catalyst activity, expressed by:

$$f_c = kc_i$$
, then

using mass balance of a catalyst around the reactor and a rearranged kinetic rate equation to determine catalyst feed factor, by:

$$f_c = kc_i = kc_o = \frac{P_o}{\tau_c C_F} = \frac{58034 \,\text{lb/h}}{(45.9 \,\text{min})(5.0537 \,\text{wt\%})} = 250 \,\text{lb/h/min/wt\%}$$

where a catalyst activity factor quantifies reactivity of a specific catalyst with respect to its size, expressed as:

$$k_c = \frac{k}{d_c^3} = \frac{7.00 \,\text{lb/lb/min/wt\%}}{(100 \,\mu)^3} \cdot \frac{\text{le06 lb}}{1 \,\text{MMlb}} = 7.00 \,\text{lb/MMlb/min/wt\%/} \mu^3$$

where μ^3 is expressed in microns.